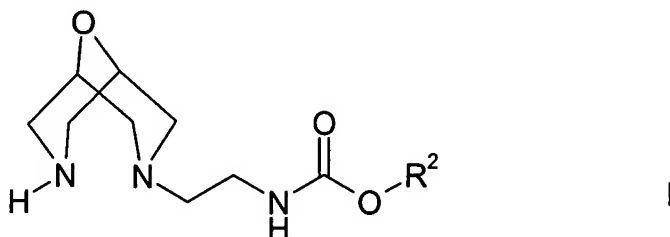


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

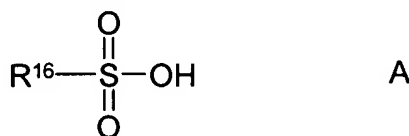
Listing of Claims:

1. (currently amended) An acid addition salt of a compound of Formula I



wherein R² represents C₁₋₆ alkyl (optionally substituted ~~and/or terminated~~ by one or more substituents selected from -OH, halo, cyano, nitro and aryl) or aryl, wherein each aryl and aryloxy group, ~~unless otherwise specified,~~ is optionally substituted.

2. (currently amended) A salt according to claim 1, wherein ~~in which~~ the acid component of the acid addition salt is represented by formula A



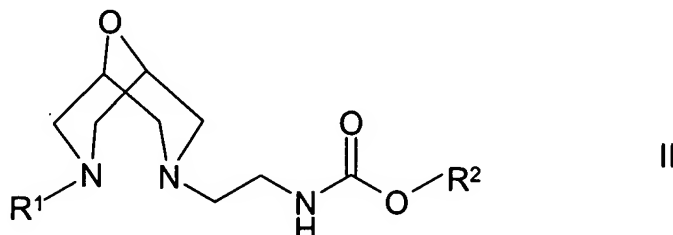
wherein R¹⁶ represents unsubstituted C₁₋₄ alkyl, C₁₋₄ perfluoroalkyl or phenyl, which latter group is optionally substituted by one or more substituents selected from C₁₋₆ alkyl, halo, nitro and C₁₋₆ alkoxy, ~~and R² is as defined above.~~

3. (currently amended) A salt according to claim 2, wherein the salt is a toluenesulfonate, benzenesulfonate, nosylate, brosylate, besylate or mesitylate salt.

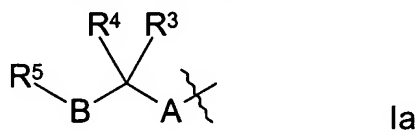
4. (currently amended) A salt according to ~~any previous~~ claim 1, wherein ~~in which~~ the salt is in solid form.

5. (currently amended) A salt according to ~~any previous claim 1, wherein the salt which is~~ [2-(9-oxa-3,7-diazabicyclo[3.3.1]non-3-yl)-ethyl]-carbamic acid *tert*-butyl ester 2,4,6-trimethylbenzenesulfonic acid.

6. (currently amended) A process for the preparation of a compound of Formula II



wherein R¹ represents a ~~structural fragment moiety~~ of formula Ia



~~in which~~ wherein A represents CH₂ and R³ represents -OH or -N(H)R⁷;

R⁴ represents H, C₁₋₆ alkyl or, together with R³, represents =O;

R⁵ represents phenyl or pyridyl, ~~both of which groups are~~ optionally substituted by one or more substituents selected from -OH, cyano, halo, nitro, C₁₋₆ alkyl (optionally terminated by -N(H)C(O)OR^{13a}), C₁₋₆ alkoxy, -N(R^{14a})R^{14b}, -C(O)R^{14c}, -C(O)OR^{14d}, -C(O)N(R^{14e})R^{14f}, -N(R^{14g})C(O)R^{14h}, -N(R¹⁴ⁱ)C(O)N(R^{14j})R^{14k}, -N(R^{14m})S(O)₂R^{13b}, -S(O)₂R^{13c} and/or -OS(O)₂R^{13d};

R⁷ represents H, C₁₋₆ alkyl, -E-aryl, -E-Het¹, -C(O)R^{9a}, -C(O)OR^{9b}, -S(O)₂R^{9c}, -[C(O)]_pN(R^{10a})R^{10b} or -C(NH)NH₂;

R^{9a} to R^{9d} independently represent, independently at each occurrence ~~when used herein~~, C₁₋₆ alkyl (optionally substituted ~~and/or terminated~~ by one or more substituents selected from halo, aryl and Het²), aryl, Het³, or R^{9a} and R^{9d} independently represent H;

R^{10a} and R^{10b} independently represent, at each occurrence ~~when used herein~~, H or C₁₋₆ alkyl (optionally substituted ~~and/or terminated~~ by one or more substituents selected from halo, aryl and Het⁴), aryl, Het⁵, or together represent C₃₋₆ alkylene, optionally interrupted by an O atom;

E represents, independently at each occurrence ~~when used herein~~, a direct bond or C₁₋₄ alkylene;

B represents -Z-, -Z-N(R¹²)-, -N(R¹²)-Z-, -Z-S(O)_n- or -Z-O- (in which latter two groups, Z is attached to the carbon atom bearing R³ and R⁴);

Z represents a direct bond or C₁₋₄ alkylene;

R¹¹ and R¹² independently represent H or C₁₋₆ alkyl;

R^{13a} to R^{13d} independently represent C₁₋₆ alkyl;

R^{14a} and R^{14b} independently represent H, C₁₋₆ alkyl or together represent C₃₋₆ alkylene, resulting in a four- to seven-membered nitrogen-containing ring;

R^{14c} to R^{14m} independently represent H or C₁₋₆ alkyl; and

n represents 0, 1 or 2;

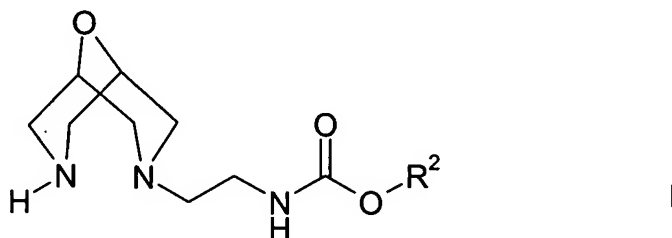
p represents 1 or 2;

Het¹ to Het⁵ independently represent, independently at each occurrence ~~when used herein~~, five- to twelve-membered heterocyclic groups containing one or more heteroatoms selected from oxygen, nitrogen and/or sulfur, which heterocyclic groups are optionally substituted by one or more substituents selected from =O, -OH, cyano, halo, nitro, C₁₋₆ alkyl, C₁₋₆ alkoxy, aryl, aryloxy, -N(R^{15a})R^{15b}, -C(O)R^{15c}, -C(O)OR^{15d}, -C(O)N(R^{15e})R^{15f}, -N(R^{15g})C(O)R^{15h} and -N(R¹⁵ⁱ)S(O)₂R^{15j};

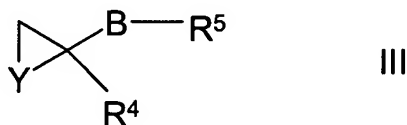
R^{15a} to R^{15j} independently represent C₁₋₆ alkyl, aryl or R^{15a} to R¹⁵ⁱ independently represent H; and

R² represents C₁₋₆ alkyl (optionally substituted ~~and/or terminated~~ by one or more substituents selected from -OH, halo, cyano, nitro and aryl) or aryl, wherein each aryl and aryloxy group, ~~unless otherwise specified~~, is optionally substituted.

wherein a salt of a compound of Formula I



~~in which~~ wherein R² is as previously defined is reacted with a compound of Formula III



wherein Y represents O or N(R⁷) and R⁴, R⁵, R⁷ and B are as hereinbefore defined,

at a temperature in the range of 0 °C to 100 °C ~~for example at elevated temperature (e.g. 60°C to reflux)~~ in the presence of a water and in the presence of a base.

7. (currently amended) A process according to claim 6, wherein ~~in which~~ the salt has been previously isolated in solid form.
8. (currently amended) A process according to ~~either claim 6 or claim 7~~ for the preparation of *tert*-butyl 2-{7-[(2*S*)-3-(4-cyanophenoxy)-2-hydroxypropyl]-9-oxa-3,7-diaza-bicyclo[3.3.1]-non-3-yl}ethylcarbamate which comprises reacting a salt of [2-(9-oxa-3,7-diazabicyclo[3.3.1]non-3-yl)-ethyl]-carbamic acid *tert*-butyl ester with 4-[(2*S*)-oxiranylmethoxy]benzonitrile at a temperature in the range of 0 °C to 100 °C in the presence of water and in the presence of a base.
9. (currently amended) A process according to ~~any one of claims~~ claim 6, wherein 7 or 8 in which the salt of Formula I is an isolated salt of [2-(9-oxa-3,7-diazabicyclo[3.3.1]non-3-yl)-ethyl]-carbamic acid *tert*-butyl ester ~~is used~~.
10. (currently amended) A process according to ~~either~~ claim 9, wherein the salt is the 2,4,6-trimethylbenzenesulfonic acid salt.